In response to the Office Action dated January 10, 2008 (Paper No.

20080104), please amend the above-identified application as follows:

## a.) Amendment to the Claims

Claims 1-10 (Cancelled).

11. (Currently Amended) A benzoyl compound represented by general formula (IA):

[wherein

nA represents an integer of 1 to 5;

R<sup>1A</sup> represents substituted or unsubstituted lower alkyl, substituted or unsubstituted lower alkoxy, substituted or unsubstituted cycloalkyl, substituted or unsubstituted lower alkoxycarbonyl, substituted or unsubstituted heterocyclic alkyl, substituted or unsubstituted aryl, CONR<sup>7</sup>R<sup>8</sup> (wherein R<sup>7</sup> and R<sup>8</sup> independently represent a hydrogen atom, substituted or unsubstituted lower alkyl, substituted or unsubstituted cycloalkyl, substituted or unsubstituted lower alkanoyl, substituted or unsubstituted aryl, a substituted or unsubstituted heterocyclic group, substituted or unsubstituted aralkyl, substituted or unsubstituted heterocyclic alkyl, or substituted or unsubstituted aroyl, or R<sup>7</sup>

and  $R^8$  form a substituted or unsubstituted heterocyclic group together with the adjacent nitrogen atom) or  $NR^9R^{10}$  (wherein  $R^9$  and  $R^{10}$  have the same meanings as the above  $R^7$  and  $R^8$ , respectively);

R<sup>2A</sup> represents substituted or unsubstituted aryl;

 $R^{3A}$  and  $R^{5A}$   $R^{3A}$ ,  $R^{4A}$  and  $R^{5A}$  each represent a hydrogen atom; and

R<sup>4A</sup> represents a hydrogen atom; and

R<sup>6A</sup> represents a hydrogen atom, halogen, lower alkyl or aryl; provided that:

when R is a hydrogen atom, and (CH<sub>2</sub>)<sub>nA</sub>R is

(a) 2 (acetoxymethyl)heptyl, 3 oxopentyl or pentyl,

R<sup>2A</sup>-is not 6-hydroxy-4-methoxy-3-methoxycarbonyl-2-pentylphenyl,

(b) 3-oxopentyl,

 $R^{\frac{2A}{1}}$  is not a group selected from the group consisting of 3-benzyloxycarbonyl 6 hydroxy 4 methoxy 2 pentylphenyl and 3 carboxy 6 hydroxy 4 methoxy 2 pentylphenyl, and

(c) n propyl,

R<sup>2A</sup>-is not 2,4 dihydroxy 6 [(4 hydroxy 2 oxopyran 6-yl)methyl]phenyl],

or a pharmaceutically acceptable salt thereof.

12. (Previously Presented) The benzoyl compound according to claim 11, wherein  $R^{2A}$  is substituted aryl having 1 to 3 substituents, or aryl, or a pharmaceutically acceptable salt thereof.

Claims 13 and 14 (Cancelled).

- 15. (Previously Presented) The benzoyl compound according to claim 11 or 12, wherein  $R^{1A}$  is  $CONR^{7}R^{8}$ , or a pharmaceutically acceptable salt thereof.
- 16. (Previously Presented) The benzoyl compound according to claim  $R^{6A}$  is lower alkyl, or a pharmaceutically acceptable salt thereof.
- 17. (Previously Presented) A pharmaceutical composition comprising, as an active ingredient, the benzoyl compound according to claim 16 or a pharmaceutically

acceptable salt of said benzoyl compound, together with a pharmaceutically acceptable carrier.

Claims 18-31 (Cancelled).

32. (Previously Presented) The benzoyl compound according to claim11, wherein

nA is 1,

 $R^{1A}$  is  $CONR^{7a}\,R^{8a}$  (wherein  $R^{7a}$  and  $R^{8a}$  independently represent lower alkyl substituted by hydroxyl or lower alkoxy), and

R<sup>6A</sup> is lower alkyl,

or a pharmaceutically acceptable salt thereof.

33. (Previously Presented) The benzoyl compound according to claim

32, wherein R<sup>2A</sup> is

(wherein  $R^{2a}$  represents substituted or unsubstituted lower alkoxy, and  $R^{2b}$  represents substituted or unsubstituted heterocyclic-alkyl oxy), or a pharmaceutically acceptable salt thereof.

34. (Previously Presented) The benzoyl compound according to claim

32, wherein 
$$\mathbb{R}^{2A}$$
 is

(wherein  $R^{2a}$  represents substituted or unsubstituted lower alkoxy, and  $R^{2b}$  represents substituted or unsubstituted heterocyclic-alkyl oxy), or a pharmaceutically acceptable salt thereof.

35. (Previously Presented) The benzoyl compound according to claim

32, wherein 
$$R^{2A}$$
 is

(wherein  $R^{2b1}$  represents substituted or unsubstituted morpholino or substituted or unsubstituted piperidino), or a pharmaceutically acceptable salt thereof.